

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111315\
 Data File : BE091161.D
 Acq On : 14 Nov 2015 3:22
 Operator : UM/NP
 Sample : G4342-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 D9KQ9

Manual Integrations
 APPROVED

mmdadoda
 11/16/2015 3:25:12 PM

Quant Time: Nov 16 11:08:19 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 14 03:28:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	4033	0.40	ng/ul	-0.07
4) Naphthalene-d8	9.56	136	20639	0.40	ng/ul	-0.05
8) Acenaphthene-d10	13.38	164	12707	0.40	ng/ul	-0.02
12) Phenanthrene-d10	16.10	188	29820m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.24	240	47761	0.40	ng/ul	0.00
22) Perylene-d12	22.13	264	42705m	0.40	ng/ul	-0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.05	96	24192	8.34	ng/uL	-0.11
6) 2-Methylnaphthalene-d10	11.09	152	6757	0.24	ng/ul	-0.03
16) Fluoranthene-d10	18.08	212	18504	0.22	ng/ul	-0.01
Target Compounds				Ovalue		
14) Phenanthrene	16.12	178	32163	0.09	ng/ul	97
17) Fluoranthene	18.11	202	62618	0.16	ng/ul	87
19) Pyrene	18.49	202	69352	0.13	ng/ul#	76
20) Benzo(a)anthracene	20.21	228	6087m	0.05	ng/ul	
21) Chrysene	20.27	228	8809	0.06	ng/ul	97
23) Benzo(b)fluoranthene	21.57	252	8097	0.06	ng/ul	72
24) Benzo(k)fluoranthene	21.60	252	7722	0.06	ng/ul#	69
25) Benzo(a)pyrene	22.03	252	7392	0.06	ng/ul#	71
26) Indeno(1,2,3-cd)pyrene	23.74	276	22070	0.04	ng/ul#	88
28) Benzo(g,h,i)perylene	24.30	276	26064	0.05	ng/ul#	85

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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